

Data Path : Z:\HPCHEM1\BNA E\DATA\BE092617\
 Data File : BE094124.D
 Acq On : 26 Sep 2017 9:46
 Operator : SJ/JU
 Sample : I5340-06RE
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleID :
 LT-TS-006

Manual Integrations
 APPROVED

Sohil
 9/27/2017 12:45:34 PM

Quant Time: Sep 26 16:22:59 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE092517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 25 17:37:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	1501	0.40	ng	0.00
7) Naphthalene-d8	10.69	136	7204	0.40	ng	0.00
13) Acenaphthene-d10	14.51	164	4800	0.40	ng	0.00
19) Phenanthrene-d10	17.24	188	18340	0.40	ng	0.00
27) Chrysene-d12	21.40	240	15729	0.40	ng	0.00
34) Perylene-d12	23.91	264	14895	0.40	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Fluorophenol	5.50	112	2371	0.42	ng	0.00
5) Phenol-d6	7.09	99	3374	0.43	ng	0.00
8) Nitrobenzene-d5	9.06	82	2819	0.40	ng	0.00
11) 2-Methylnaphthalene-d10	12.27	152	4923	0.44	ng	0.00
14) 2,4,6-Tribromophenol	16.00	330	1019	0.37	ng	0.00
15) 2-Fluorobiphenyl	13.13	172	6427	0.41	ng	0.00
25) Fluoranthene-d10	19.27	212	71186	0.40	ng	0.00
29) Terphenyl-d14	19.85	244	11061	0.39	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
9) Naphthalene	10.74	128	3090	0.038	ng	# 92
12) 2-Methylnaphthalene	12.35	142	498	0.038	ng	# 85
16) Acenaphthylene	14.23	152	5238	0.218	ng	99
17) Acenaphthene	14.57	154	649	0.043	ng	# 91
18) Fluorene	15.56	166	1542	0.076	ng	90
23) Phenanthrene	17.27	178	45646	0.973	ng	98
24) Anthracene	17.37	178	6990	0.142	ng	# 92
26) Fluoranthene	19.30	202	101640	1.933	ng	99
28) Pyrene	19.66	202	100657	2.020	ng	# 96
30) Benzo(a)anthracene	21.38	228	47371	0.894	ng	97
31) Chrysene	21.44	228	57588	1.127	ng	97
32) Bis(2-ethylhexyl)phthalate	21.29	149	42086	0.059	ng	99
33) Indeno(1,2,3-cd)pyrene	26.55	276	33631	0.615	ng	96
35) Benzo(b)fluoranthene	23.14	252	73183	1.362	ng	97
36) Benzo(k)fluoranthene	23.18	252	24263m	0.472	ng	
37) Benzo(a)pyrene	23.79	252	48855	1.006	ng	99
38) Dibenzo(a,h)anthracene	26.56	278	8508	0.185	ng	# 77
39) Benzo(a,h,i)perylene	27.38	276	34556	0.759	ng	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\HPCHEM1\BNA E\DATA\BE092617\
Data File : BE094124.D
Acq On : 26 Sep 2017 9:46
Operator : SJ/JU
Sample : I5340-06RE
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_E
ClientSampled :
LT-TS-006

Manual Integrations
APPROVED

Sohil
9/27/2017 12:45:34 PM

Quant Time: Sep 26 16:22:59 2017
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE092517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Sep 25 17:37:08 2017
Response via : Initial Calibration

