

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE032319\
 Data File : BE099112.D
 Acq On : 23 Mar 2019 15:09
 Operator : JU/SJ
 Sample : K2014-07
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PST-026-SW040-031919

Manual Integrations
 APPROVED

Sohil
 3/25/2019 12:36:59 PM

Quant Time: Mar 24 03:18:33 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE032219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 23 04:40:11 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	3226	0.40	ng	0.00
7) Naphthalene-d8	10.63	136	12457	0.40	ng	-0.01
13) Acenaphthene-d10	14.48	164	8009	0.40	ng	0.00
19) Phenanthrene-d10	17.21	188	22985	0.40	ng	0.00
27) Chrysene-d12	21.39	240	34068	0.40	ng	0.00
34) Perylene-d12	23.90	264	36246	0.40	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.41	112	4013	0.42	ng	0.00
5) Phenol-d6	6.99	99	5637	0.41	ng	0.00
8) Nitrobenzene-d5	8.99	82	3496	0.64	ng	0.00
11) 2-Methylnaphthalene-d10	12.22	152	8242	0.37	ng	-0.02
14) 2,4,6-Tribromophenol	15.96	330	1257	0.53	ng	0.00
15) 2-Fluorobiphenyl	13.10	172	12900	0.40	ng	-0.02
25) Fluoranthene-d10	19.24	212	119489	0.39	ng	0.00
29) Terphenyl-d14	19.83	244	22890	0.32	ng	0.00

Target Compounds				Qvalue		
9) Naphthalene	10.68	128	3526	0.024	ng	# 96
16) Acenaphthylene	14.20	152	1785	0.042	ng	97
18) Fluorene	15.51	166	933	0.025	ng	# 89
23) Phenanthrene	17.25	178	23249	0.341	ng	99
24) Anthracene	17.35	178	1998	0.031	ng	# 89
26) Fluoranthene	19.27	202	38865	0.405	ng	99
28) Pyrene	19.63	202	41414	0.358	ng	96
30) Benzo(a)anthracene	21.37	228	20441	0.170	ng	99
31) Chrysene	21.43	228	23016	0.192	ng	98
32) Bis(2-ethylhexyl)phthalate	21.29	149	18779	0.121	ng	99
33) Indeno(1,2,3-cd)pyrene	26.55	276	11187	0.087	ng	97
35) Benzo(b)fluoranthene	23.12	252	23908	0.191	ng	94
36) Benzo(k)fluoranthene	23.17	252	10109m	0.082	ng	
37) Benzo(a)pyrene	23.78	252	17713	0.153	ng	99
38) Dibenzo(a,h)anthracene	26.57	278	2891	0.026	ng	# 77
39) Benzo(g,h,i)perylene	27.37	276	12207	0.105	ng	99

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

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