

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN081319\
 Data File : BN007154.D
 Acq On : 14 Aug 2019 05:45
 Operator : HP/JU
 Sample : K4173-20
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 EXT-015-SW088-080119

Quant Time: Aug 14 09:21:34 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN081019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 13 12:12:02 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.43	152	9423	0.40	ng	0.00
6) Naphthalene-d8	10.19	136	36996	0.40	ng	0.00
12) Acenaphthene-d10	14.08	164	21585	0.40	ng	0.00
18) Phenanthrene-d10	16.82	188	47765	0.40	ng	0.00
26) Chrysene-d12	21.05	240	46442	0.40	ng	-0.01
32) Perylene-d12	23.16	264	56563	0.40	ng	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 2-Fluorophenol	5.06	112	6425	0.28	ng	0.00
4) Phenol-d6	6.62	99	8533	0.30	ng	0.00
7) Nitrobenzene-d5	8.56	82	6903	0.23	ng	0.00
10) 2-Methylnaphthalene-d10	11.79	152	15072	0.22	ng	0.00
13) 2,4,6-Tribromophenol	15.57	330	2511	0.22	ng	0.00
14) 2-Fluorobiphenyl	12.71	172	3627	0.12	ng	0.00
24) Fluoranthene-d10	18.87	212	39425	0.23	ng	0.00
28) Terphenyl-d14	19.49	244	28508	0.23	ng	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
8) Naphthalene	10.24	128	32422	0.313	ng	99
11) 2-Methylnaphthalene	11.87	142	37271	0.507	ng	98
15) Acenaphthylene	13.79	152	331112	2.935	ng	99
16) Acenaphthene	14.13	154	45676	0.632	ng	94
17) Fluorene	15.13	166	114569	1.060	ng	88
22) Phenanthrene	16.87	178	1915265	13.380	ng	100
23) Anthracene	16.95	178	379290	2.900	ng	99
25) Fluoranthene	18.90	202	2010067	10.977	ng	98
27) Pyrene	19.26	202	2850922	17.505	ng	96
29) Benzo(a)anthracene	21.02	228	1164617	6.857	ng	93
30) Chrysene	21.08	228	1454851	8.817	ng	96
31) Indeno(1,2,3-cd)pyrene	25.24	276	657596	3.616	ng	99
33) Benzo(b)fluoranthene	22.54	252	1320233	6.807	ng	99
34) Benzo(k)fluoranthene	22.58	252	349144m	1.823	ng	
35) Benzo(a)pyrene	23.07	252	1052756	5.987	ng	99
36) Dibenzo(a,h)anthracene	25.25	278	191653	1.092	ng	# 88
37) Benzo(g,h,i)perylene	25.87	276	746353	4.136	ng	99

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

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