

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE042619\
 Data File : BE099385.D
 Acq On : 26 Apr 2019 19:19
 Operator : JU/SJ
 Sample : K2569-06
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PST-003-SW024-042319

Manual Integrations
 APPROVED

Sohil
 4/29/2019 1:28:50 PM

Quant Time: Apr 27 06:50:35 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE041819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 19 06:51:38 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.82	152	4182	0.40	ng	0.00
7) Naphthalene-d8	10.61	136	16136	0.40	ng	0.01
13) Acenaphthene-d10	14.46	164	10969	0.40	ng	0.00
19) Phenanthrene-d10	17.19	188	29241	0.40	ng	0.00
27) Chrysene-d12	21.37	240	32236	0.40	ng	0.00
34) Perylene-d12	23.87	264	37094	0.40	ng	0.01

System Monitoring Compounds						
4) 2-Fluorophenol	5.40	112	3178	0.31	ng	0.01
5) Phenol-d6	6.98	99	4923	0.34	ng	0.00
8) Nitrobenzene-d5	8.98	82	4081	0.33	ng	0.01
11) 2-Methylnaphthalene-d10	12.21	152	7503	0.26	ng	0.01
14) 2,4,6-Tribromophenol	15.94	330	1712	0.47	ng	0.00
15) 2-Fluorobiphenyl	13.09	172	11080	0.26	ng	0.00
25) Fluoranthene-d10	19.22	212	88416	0.25	ng	0.00
29) Terphenyl-d14	19.82	244	18857	0.30	ng	0.00

Target Compounds				Qvalue		
9) Naphthalene	10.65	128	43797	0.249	ng	100
12) 2-Methylnaphthalene	12.28	142	6253	0.209	ng	97
16) Acenaphthylene	14.18	152	1549	0.031	ng	# 83
17) Acenaphthene	14.53	154	6852	0.224	ng	99
18) Fluorene	15.50	166	7193	0.164	ng	98
23) Phenanthrene	17.24	178	143672	1.855	ng	99
24) Anthracene	17.32	178	20123	0.292	ng	# 90
26) Fluoranthene	19.25	202	187238	1.807	ng	98
28) Pyrene	19.62	202	160477	1.690	ng	# 95
30) Benzo(a)anthracene	21.36	228	95748	1.010	ng	98
31) Chrysene	21.40	228	78340	0.777	ng	96
32) Bis(2-ethylhexyl)phthalate	21.27	149	16482	0.191	ng	# 100
33) Indeno(1,2,3-cd)pyrene	26.52	276	36088	0.362	ng	# 92
35) Benzo(b)fluoranthene	23.10	252	94394	0.831	ng	95
36) Benzo(k)fluoranthene	23.15	252	31132m	0.275	ng	
37) Benzo(a)pyrene	23.75	252	67506	0.663	ng	96
38) Dibenzo(a,h)anthracene	26.54	278	8830	0.085	ng	# 75
39) Benzo(a,h,i)perylene	27.34	276	33048	0.313	ng	97

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

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