

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE052219\
 Data File : BE099717.D
 Acq On : 22 May 2019 13:20
 Operator : JU/SJ
 Sample : K2966-06
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 FESB-18(2.5-3.0)

Manual Integrations
 APPROVED

mohammad
 5/24/2019 8:47:56 AM

Quant Time: May 23 08:08:42 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE052019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 20 19:29:40 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.82	152	1670	0.40	ng	0.00
7) Naphthalene-d8	10.61	136	6061	0.40	ng	0.00
13) Acenaphthene-d10	14.49	164	4715	0.40	ng	0.00
19) Phenanthrene-d10	17.23	188	13828	0.40	ng	0.00
27) Chrysene-d12	21.43	240	24068	0.40	ng	0.01
34) Perylene-d12	23.96	264	25107	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.37	112	1513	0.36	ng	0.00
5) Phenol-d6	6.98	99	2181	0.39	ng	0.00
8) Nitrobenzene-d5	8.98	82	1785	0.34	ng	0.00
11) 2-Methylnaphthalene-d10	12.21	152	2645	0.25	ng	-0.02
14) 2,4,6-Tribromophenol	15.97	330	961	0.40	ng	-0.01
15) 2-Fluorobiphenyl	13.10	172	3734	0.21	ng	0.00
25) Fluoranthene-d10	19.27	212	30081	0.16	ng	0.00
29) Terphenyl-d14	19.86	244	8018m	0.19	ng	-0.01
Target Compounds						Qvalue
9) Naphthalene	10.67	128	72759	1.123	ng	99
12) 2-Methylnaphthalene	12.30	142	8740	0.827	ng	96
16) Acenaphthylene	14.21	152	65363	2.974	ng	100
17) Acenaphthene	14.54	154	37386	3.031	ng	98
18) Fluorene	15.51	166	79670	4.385	ng	99
22) Pentachlorophenol	16.91	266	820	0.503	ng	# 36
23) Phenanthrene	17.27	178	1437749	42.549	ng	99
24) Anthracene	17.36	178	240791	7.858	ng	# 95
26) Fluoranthene	19.30	202	3468118	64.921	ng	97
28) Pvrene	19.66	202	3302077	54.607	ng	# 91
30) Benzo(a)anthracene	21.40	228	2197881	32.137	ng	95
31) Chrysene	21.46	228	2124172	29.050	ng	96
32) Bis(2-ethylhexyl)phthalate	21.32	149	249056	3.160	ng	100
33) Indeno(1,2,3-cd)pyrene	26.67	276	1365475	15.468	ng	# 96
35) Benzo(b)fluoranthene	23.19	252	2654910	38.589	ng	98
36) Benzo(k)fluoranthene	23.23	252	942721m	13.145	ng	
37) Benzo(a)pvrene	23.86	252	2012979m	30.405	ng	
38) Dibenzo(a,h)anthracene	26.68	278	360569	5.187	ng	93
39) Benzo(g,h,i)perylene	27.52	276	1318869	18.786	ng	98

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

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