

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092517\
 Data File : BF098800.D
 Acq On : 25 Sep 2017 23:27
 Operator : SJ/JU
 Sample : I5380-01 50X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-04-092217

Manual Integrations
 APPROVED

Sohil
 9/26/2017 4:58:42 PM

Quant Time: Sep 26 07:50:54 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 23 13:50:58 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.91	152	150236	20.00	ng	-0.01
21) Naphthalene-d8	8.19	136	645493	20.00	ng	-0.01
38) Acenaphthene-d10	9.95	164	307984	20.00	ng	-0.01
63) Phenanthrene-d10	11.44	188	555585	20.00	ng	0.00
75) Chrysene-d12	14.08	240	338654	20.00	ng	0.00
86) Perylene-d12	15.57	264	276466	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.52	112	13855	1.47	ng	-0.02
7) Phenol-d6	6.52	99	19895	1.71	ng	-0.02
23) Nitrobenzene-d5	7.47	82	11266	1.12	ng	-0.02
41) 2,4,6-Tribromophenol	10.73	330	5515	2.19	ng	-0.02
44) 2-Fluorobiphenyl	9.26	172	33215	1.80	ng	-0.02
78) Terphenyl-d14	13.02	244	36609	2.46	ng	-0.02

Target Compounds

						Qvalue
31) Naphthalene	8.22	128	133817	4.414	ng	98
37) 2-Methylnaphthalene	8.90	142	40213	2.040	ng	99
51) Acenaphthene	9.98	154	120176	6.236	ng	98
54) Dibenzofuran	10.15	168	117587	4.583	ng	98
57) Fluorene	10.50	166	172362	8.357	ng	99
70) Phenanthrene	11.47	178	1429588	48.071	ng	99
71) Anthracene	11.52	178	506160	16.634	ng	99
72) Carbazole	11.66	167	125500	4.212	ng	99
74) Fluoranthene	12.66	202	1968900	65.382	ng	99
77) Pyrene	12.89	202	1672560	64.769	ng	100
80) Benzo(a)anthracene	14.07	228	694010	33.844	ng	98
82) Chrysene	14.10	228	550768	28.211	ng	97
85) Indeno(1,2,3-cd)pyrene	17.07	276	282306	18.077	ng	96
87) Benzo(b)fluoranthene	15.13	252	552842m	32.523	ng	
88) Benzo(k)fluoranthene	15.16	252	219449m	13.071	ng	
89) Benzo(a)pyrene	15.50	252	416802	26.713	ng	# 94
90) Dibenzo(a,h)anthracene	17.08	278	65450	5.241	ng	94
91) Benzo(g,h,i)perylene	17.53	276	262635	21.278	ng	97

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

