

Data Path : Z:\HPCHEM1\BNA F\DATA\BF081617\
 Data File : BF097772.D
 Acq On : 17 Aug 2017 7:18
 Operator : SJ/JU
 Sample : I4791-06
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-22-COMP

Manual Integrations
 APPROVED

Sohil
 8/17/2017 5:58:11 PM

Quant Time: Aug 17 14:52:12 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 15 13:25:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.78	152	150897	20.00	ng	0.00
21) Naphthalene-d8	8.06	136	572578	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	229992	20.00	ng	0.00
63) Phenanthrene-d10	11.30	188	339913	20.00	ng	0.00
75) Chrysene-d12	13.93	240	288622	20.00	ng	0.00
86) Perylene-d12	15.36	264	246046	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	1317676	132.82	ng	0.01
7) Phenol-d6	6.42	99	1779485	132.02	ng	0.00
23) Nitrobenzene-d5	7.35	82	1073005	101.80	ng	0.00
41) 2,4,6-Tribromophenol	10.60	330	244425	122.48	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	1477848	94.41	ng	0.00
78) Terphenyl-d14	12.89	244	852297	61.58	ng	0.00
Target Compounds						
10) Phenol	6.43	94	50851	3.226	ng	92
49) Dimethylphthalate	9.53	163	208954	12.429	ng	99
74) Fluoranthene	12.50	202	57962	3.066	ng	96
77) Pyrene	12.73	202	68157	2.887	ng	97
80) Benzo(a)anthracene	13.92	228	41175m	2.380	ng	
82) Chrysene	13.96	228	42594	2.475	ng	95
85) Indeno(1,2,3-cd)pyrene	16.75	276	30173m	2.384	ng	
87) Benzo(b)fluoranthene	14.95	252	65940m	4.252	ng	
89) Benzo(a)pyrene	15.30	252	50010	3.642	ng	98
91) Benzo(g,h,i)perylene	17.17	276	37725	3.235	ng	95

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

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