

Data Path : Z:\HPCHEM1\BNA F\DATA\BF120717\
 Data File : BF101201.D
 Acq On : 7 Dec 2017 13:27
 Operator : SJ/JU
 Sample : I6615-01 2X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB-2-OW(0-10)

Manual Integrations
 APPROVED

Sohil
 12/8/2017 11:12:00 AM

Quant Time: Dec 07 16:49:30 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 04 13:07:49 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	53958	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	210621	20.00	ng	-0.01
38) Acenaphthene-d10	9.88	164	98007	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	177795	20.00	ng	0.00
75) Chrysene-d12	14.02	240	115267	20.00	ng	0.00
86) Perylene-d12	15.50	264	114516	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	183603	56.23	ng	0.00
7) Phenol-d6	6.47	99	223716	56.43	ng	0.00
23) Nitrobenzene-d5	7.40	82	132993	37.67	ng	-0.01
41) 2,4,6-Tribromophenol	10.67	330	60707	51.56	ng	0.00
44) 2-Fluorobiphenyl	9.20	172	292277	40.80	ng	-0.01
78) Terphenyl-d14	12.96	244	235024	44.60	ng	0.00
Target Compounds						
10) Phenol	6.49	94	9030	2.048	ng	Qvalue # 69
70) Phenanthrene	11.39	178	32793	3.349	ng	99
74) Fluoranthene	12.59	202	45985	4.237	ng	94
77) Pyrene	12.82	202	44132	5.549	ng	98
80) Benzo(a)anthracene	14.01	228	18392	2.527	ng	94
82) Chrysene	14.04	228	18411	2.724	ng	98
87) Benzo(b)fluoranthene	15.06	252	21358m	3.114	ng	
89) Benzo(a)pyrene	15.43	252	16626	2.666	ng	96
91) Benzo(g,h,i)perylene	17.43	276	10999	2.123	ng	Qvalue # 90

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

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