

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110217\
 Data File : BF100103.D
 Acq On : 2 Nov 2017 23:33
 Operator : SJ/JU
 Sample : I6216-01
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EPE-109-3(0-5)

Manual Integrations
 APPROVED

Sohil
 11/3/2017 11:20:30 AM

Quant Time: Nov 03 04:03:28 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 01 18:13:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	94678	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	377189	20.00	ng	-0.01
38) Acenaphthene-d10	9.80	164	166800	20.00	ng	-0.01
63) Phenanthrene-d10	11.30	188	263611	20.00	ng	-0.01
75) Chrysene-d12	13.94	240	164228	20.00	ng	-0.03
86) Perylene-d12	15.41	264	165276	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	428924	71.12	ng	0.00
7) Phenol-d6	6.42	99	516563	72.24	ng	-0.01
23) Nitrobenzene-d5	7.34	82	281849	48.11	ng	-0.01
41) 2,4,6-Tribromophenol	10.60	330	106135	69.82	ng	-0.01
44) 2-Fluorobiphenyl	9.12	172	526938	48.74	ng	-0.01
78) Terphenyl-d14	12.88	244	359534	47.84	ng	-0.02
Target Compounds						
10) Phenol	6.43	94	20276	2.583	ng	81
49) Dimethylphthalate	9.52	163	113396	8.585	ng	100
70) Phenanthrene	11.32	178	42001	2.823	ng	98
74) Fluoranthene	12.52	202	230655	14.919	ng	97
77) Pyrene	12.74	202	222564	17.823	ng	98
80) Benzo(a)anthracene	13.94	228	117360	11.273	ng	99
82) Chrysene	13.97	228	82218	8.400	ng	97
85) Indeno(1,2,3-cd)pyrene	16.88	276	54543	6.070	ng	97
87) Benzo(b)fluoranthene	14.99	252	135809m	13.013	ng	
88) Benzo(k)fluoranthene	15.01	252	47667m	4.844	ng	
89) Benzo(a)pyrene	15.35	252	102236	10.905	ng	97
91) Benzo(a,h,i)perylene	17.33	276	47212	5.793	ng	92

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

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