

Data Path : Z:\HPCHEM1\BNA F\DATA\BF102817\
 Data File : BF099928.D
 Acq On : 28 Oct 2017 20:18
 Operator : SJ/JU
 Sample : I6167-01
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S13-SP2C

Manual Integrations
 APPROVED

Sohil
 10/30/2017 6:20:36 PM

Quant Time: Oct 30 04:26:42 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 27 15:58:57 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	100826	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	412233	20.00	ng	-0.01
38) Acenaphthene-d10	9.82	164	178525	20.00	ng	-0.02
63) Phenanthrene-d10	11.32	188	261153	20.00	ng	-0.02
75) Chrysene-d12	13.97	240	178099	20.00	ng	-0.02
86) Perylene-d12	15.43	264	181177	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	503788	78.45	ng	0.00
7) Phenol-d6	6.43	99	612514	80.44	ng	0.00
23) Nitrobenzene-d5	7.36	82	353130	55.15	ng	-0.01
41) 2,4,6-Tribromophenol	10.62	330	101548	62.42	ng	-0.02
44) 2-Fluorobiphenyl	9.14	172	502734	43.45	ng	-0.02
78) Terphenyl-d14	12.90	244	306936	37.66	ng	-0.02
Target Compounds						
10) Phenol	6.45	94	16999	2.033	ng	Qvalue 82
49) Dimethylphthalate	9.54	163	124029	8.773	ng	99
70) Phenanthrene	11.34	178	112759	7.651	ng	97
74) Fluoranthene	12.53	202	128748	8.406	ng	98
77) Pyrene	12.76	202	116539	8.605	ng	99
80) Benzo(a)anthracene	13.96	228	73783	6.535	ng	97
82) Chrysene	13.99	228	65201	6.143	ng	97
85) Indeno(1,2,3-cd)pyrene	16.91	276	40537	4.160	ng	95
87) Benzo(b)fluoranthene	15.01	252	92997m	8.129	ng	
88) Benzo(k)fluoranthene	15.03	252	36808m	3.412	ng	
89) Benzo(a)pyrene	15.37	252	65045	6.329	ng	# 89
91) Benzo(a,h,i)perylene	17.36	276	38422	4.301	ng	# 90

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

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