

Data Path : Z:\HPCHEM1\BNA F\DATA\BF102817\
 Data File : BF099919.D
 Acq On : 28 Oct 2017 16:07
 Operator : SJ/JU
 Sample : I6077-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FILL-PILE-1-COMP-1

Manual Integrations
 APPROVED

Sohil
 10/30/2017 6:20:21 PM

Quant Time: Oct 30 02:57:34 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	91168	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	381851	20.00	ng	-0.01
38) Acenaphthene-d10	9.82	164	170245	20.00	ng	-0.02
63) Phenanthrene-d10	11.32	188	269051	20.00	ng	-0.02
75) Chrysene-d12	13.96	240	160128	20.00	ng	-0.03
86) Perylene-d12	15.43	264	171067	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	597585	102.91	ng	0.00
7) Phenol-d6	6.43	99	717532	104.21	ng	0.00
23) Nitrobenzene-d5	7.36	82	413642	69.74	ng	-0.01
41) 2,4,6-Tribromophenol	10.62	330	130842	84.33	ng	-0.02
44) 2-Fluorobiphenyl	9.14	172	742694	67.31	ng	-0.02
78) Terphenyl-d14	12.90	244	475344	64.87	ng	-0.02
Target Compounds						
10) Phenol	6.45	94	17322	2.291	ng	85
49) Dimethylphthalate	9.54	163	110898	8.226	ng	98
70) Phenanthrene	11.34	178	70805	4.663	ng	98
74) Fluoranthene	12.53	202	104745	6.638	ng	96
77) Pyrene	12.76	202	95968	7.882	ng	98
80) Benzo(a)anthracene	13.95	228	40500	3.990	ng	95
82) Chrysene	13.99	228	38638	4.049	ng	98
85) Indeno(1,2,3-cd)pyrene	16.91	276	24375	2.782	ng	98
87) Benzo(b)fluoranthene	15.00	252	49359m	4.569	ng	
89) Benzo(a)pyrene	15.37	252	35150	3.622	ng	98
91) Benzo(a,h,i)perylene	17.36	276	22769	2.699	ng	97

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

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