

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090517\
 Data File : BF098272.D
 Acq On : 5 Sep 2017 15:04
 Operator : SJ/JU
 Sample : I5073-08
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-3-4-COMP

Manual Integrations
 APPROVED

Sohil
 9/6/2017 5:04:19 PM

Quant Time: Sep 06 01:35:13 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 05 17:28:20 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	106337	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	412657	20.00	ng	0.00
38) Acenaphthene-d10	9.73	164	180917	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	354332	20.00	ng	0.00
75) Chrysene-d12	13.85	240	239026	20.00	ng	-0.01
86) Perylene-d12	15.26	264	197536	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.31	112	774846	110.84	ng	0.00
7) Phenol-d6	6.35	99	1084258	114.15	ng	0.00
23) Nitrobenzene-d5	7.27	82	645923	85.03	ng	0.00
41) 2,4,6-Tribromophenol	10.53	330	210015	133.79	ng	0.00
44) 2-Fluorobiphenyl	9.06	172	947131	76.92	ng	0.00
78) Terphenyl-d14	12.80	244	845705	73.78	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.46	163	126839	9.591	ng	98
70) Phenanthrene	11.24	178	44811	2.373	ng	95
74) Fluoranthene	12.43	202	89352	4.534	ng	98
77) Pyrene	12.66	202	97061	4.965	ng	97
80) Benzo(a)anthracene	13.84	228	55708	3.889	ng	96
82) Chrysene	13.87	228	46063m	3.232	ng	
85) Indeno(1,2,3-cd)pyrene	16.62	276	25517	2.434	ng	96
87) Benzo(b)fluoranthene	14.86	252	61806m	4.964	ng	
89) Benzo(a)pyrene	15.20	252	39242	3.560	ng	98
91) Benzo(g,h,i)perylene	17.03	276	33586	3.587	ng	92

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

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