

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070218\
 Data File : BF107037.D
 Acq On : 2 Jul 2018 11:56
 Operator : JU/SJ
 Sample : J3749-06
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ZER-BOLLARD-1-2

Manual Integrations
 APPROVED

Sohil
 7/3/2018 5:29:42 PM

Quant Time: Jul 02 14:01:04 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 29 14:38:29 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	61108	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	230246	20.00	ng	-0.01
38) Acenaphthene-d10	10.00	164	104012	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	195772	20.00	ng	0.00
75) Chrysene-d12	14.15	240	184260	20.00	ng	-0.01
86) Perylene-d12	15.71	264	130641	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	70731	19.60	ng	0.00
7) Phenol-d6	6.60	99	116273	25.80	ng	-0.01
23) Nitrobenzene-d5	7.52	82	70470	17.01	ng	-0.02
41) 2,4,6-Tribromophenol	10.80	330	12280	10.19	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	149054	11.50	ng	-0.02
78) Terphenyl-d14	13.08	244	169719	22.31	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.71	163	29548	3.788	ng	Qvalue # 98
70) Phenanthrene	11.52	178	60382	6.395	ng	98
71) Anthracene	11.57	178	24407	2.532	ng	100
74) Fluoranthene	12.71	202	254633	24.466	ng	95
77) Pyrene	12.95	202	257570	21.198	ng	98
80) Benzo(a)anthracene	14.14	228	129929	11.570	ng	97
82) Chrysene	14.18	228	123026	11.908	ng	97
85) Indeno(1,2,3-cd)pyrene	17.30	276	37620	4.291	ng	# 87
87) Benzo(b)fluoranthene	15.25	252	115356m	14.215	ng	
88) Benzo(k)fluoranthene	15.28	252	43768m	5.663	ng	
89) Benzo(a)pyrene	15.64	252	71009	9.843	ng	# 92
91) Benzo(a,h,i)perylene	17.79	276	37854	6.334	ng	# 88

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

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