

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030519\
 Data File : BF112878.D
 Acq On : 6 Mar 2019 3:24
 Operator : JU/SJ
 Sample : K1705-16
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PST-028-B101-022719

Manual Integrations
 APPROVED

Sohil
 3/6/2019 12:24:52 PM

Quant Time: Mar 06 04:46:01 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 01 15:40:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	159379	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	577240	20.00	ng	0.00
39) Acenaphthene-d10	9.77	164	249489	20.00	ng	0.00
64) Phenanthrene-d10	11.25	188	477079	20.00	ng	0.00
76) Chrysene-d12	13.87	240	454576	20.00	ng	0.00
87) Perylene-d12	15.29	264	305564	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	900391	87.88	ng	0.00
7) Phenol-d6	6.37	99	1180355	91.71	ng	0.00
23) Nitrobenzene-d5	7.30	82	696174	72.17	ng	0.00
42) 2,4,6-Tribromophenol	10.56	330	301294	113.75	ng	0.00
45) 2-Fluorobiphenyl	9.10	172	1130849	72.17	ng	0.00
79) Terphenyl-d14	12.83	244	1073325	45.77	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.49	163	94476	5.136	ng	99
71) Phenanthrene	11.27	178	82181	3.462	ng	96
75) Fluoranthene	12.46	202	115625	4.547	ng	95
78) Pyrene	12.68	202	109418	2.855	ng	98
88) Benzo(b)fluoranthene	14.89	252	46582m	2.357	ng	

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

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