

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030519\
 Data File : BM019058.D
 Acq On : 06 Mar 2019 05:28
 Operator : JU/SJ
 Sample : K1704-04
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PST-027-B114-022719

Quant Time: Mar 06 06:50:30 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM021119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 05 03:36:13 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	260553	20.00	ng	0.00
21) Naphthalene-d8	10.46	136	1020273	20.00	ng	-0.01
39) Acenaphthene-d10	14.33	164	555742	20.00	ng	0.00
64) Phenanthrene-d10	17.08	188	1236681	20.00	ng	0.00
76) Chrysene-d12	21.27	240	1253828	20.00	ng	-0.01
87) Perylene-d12	23.52	264	1191356	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.27	112	1905676	119.84	ng	0.00
7) Phenol-d6	6.85	99	2654894	118.51	ng	0.00
23) Nitrobenzene-d5	8.84	82	1661917	75.32	ng	-0.01
42) 2,4,6-Tribromophenol	15.83	330	980782	122.43	ng	0.00
45) 2-Fluorobiphenyl	12.94	172	3003737	71.75	ng	-0.01
79) Terphenyl-d14	19.72	244	4300377	70.75	ng	-0.01
Target Compounds						
10) Phenol	6.88	94	203267	8.623	ng	100
50) Dimethylphthalate	13.79	163	165127	3.565	ng	100
71) Phenanthrene	17.12	178	241961	3.641	ng	100
75) Fluoranthene	19.14	202	370926	4.832	ng	99
78) Pyrene	19.51	202	338790	4.652	ng	99
81) Benzo(a)anthracene	21.26	228	189203	2.589	ng	97
83) Chrysene	21.32	228	166899	2.382	ng	98
88) Benzo(b)fluoranthene	22.84	252	185429	2.720	ng	# 94
90) Benzo(a)pyrene	23.42	252	129257	2.002	ng	# 95

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

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