

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP062220\
 Data File : BP002486.D
 Acq On : 22 Jun 2020 18:40
 Operator : CG/JU
 Sample : L3079-08
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JFK-SB-01-0.0-72.0

Quant Time: Jun 23 01:05:21 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP060520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 05 15:08:53 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	132880	20.00	ng	0.00
21) Naphthalene-d8	10.36	136	534044	20.00	ng	0.00
39) Acenaphthene-d10	14.23	164	329038	20.00	ng	0.00
64) Phenanthrene-d10	16.99	188	721634	20.00	ng	0.00
76) Chrysene-d12	21.10	240	738902	20.00	ng	-0.01
86) Perylene-d12	23.30	264	797228	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.21	112	869341	121.06	ng	0.00
7) Phenol-d6	6.78	99	1159727	121.30	ng	0.00
23) Nitrobenzene-d5	8.73	82	797055	89.13	ng	-0.01
42) 2,4,6-Tribromophenol	15.73	330	451916	143.45	ng	0.00
45) 2-Fluorobiphenyl	12.85	172	1756979	89.78	ng	0.00
79) Terphenyl-d14	19.58	244	2675334	96.98	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.70	163	231151	10.253	ng	99
71) Phenanthrene	17.03	178	79936	2.368	ng	98
75) Fluoranthene	19.02	202	167018	4.145	ng	98
78) Pyrene	19.37	202	163017	3.602	ng	98
81) Benzo(a)anthracene	21.08	228	90813	2.153	ng	99
83) Chrysene	21.13	228	81458	2.039	ng	97
88) Benzo(b)fluoranthene	22.64	252	100660	2.341	ng	98

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

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