

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP062620\
 Data File : BP002555.D
 Acq On : 26 Jun 2020 17:22
 Operator : CG/JU
 Sample : L3126-11
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-4-S

Quant Time: Jun 26 18:19:16 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP060520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 24 13:54:06 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	218092	20.00	ng	0.00
21) Naphthalene-d8	10.35	136	875793	20.00	ng	0.00
39) Acenaphthene-d10	14.23	164	535878	20.00	ng	0.00
64) Phenanthrene-d10	16.99	188	1146300	20.00	ng	0.00
76) Chrysene-d12	21.10	240	891904	20.00	ng	0.00
86) Perylene-d12	23.30	264	774536	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.20	112	1218965	103.42	ng	0.00
7) Phenol-d6	6.78	99	1688101	107.58	ng	0.00
23) Nitrobenzene-d5	8.73	82	1191806	81.27	ng	0.00
42) 2,4,6-Tribromophenol	15.73	330	651434	126.97	ng	0.00
45) 2-Fluorobiphenyl	12.85	172	2480037	77.82	ng	0.00
79) Terphenyl-d14	19.58	244	3185783	95.11	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.69	163	267358	7.281	ng	99
71) Phenanthrene	17.03	178	214030	3.991	ng	99
75) Fluoranthene	19.02	202	344304	5.379	ng	98
78) Pyrene	19.37	202	300376	5.498	ng	100
81) Benzo(a)anthracene	21.08	228	128420	2.522	ng	98
83) Chrysene	21.13	228	121567	2.520	ng	98
88) Benzo(b)fluoranthene	22.64	252	135098	3.234	ng	# 89
90) Benzo(a)pyrene	23.20	252	91643	2.341	ng	# 95

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

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