

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062524\
 Data File : BN032262.D
 Acq On : 26 Jun 2024 05:19
 Operator : MA/JU
 Sample : P3056-02
 Misc :
 ALS Vial : 66 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BP-VPB-196-GW-665-667

Quant Time: Jun 26 05:50:48 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN061024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 19 11:30:34 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.632	152	8194	0.400	ng	-0.01
7) Naphthalene-d8	10.410	136	24493	0.400	ng	-0.01
13) Acenaphthene-d10	14.274	164	14375	0.400	ng	-0.01
19) Phenanthrene-d10	17.029	188	30193	0.400	ng	0.00
29) Chrysene-d12	21.229	240	28757	0.400	ng	# 0.00
35) Perylene-d12	23.449	264	31565	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.241	112	3514	0.151	ng	0.00
5) Phenol-d6	6.823	99	2278	0.075	ng	0.00
8) Nitrobenzene-d5	8.798	82	3395	0.167	ng	0.00
11) 2-Methylnaphthalene-d10	12.001	152	6390	0.164	ng	-0.02
14) 2,4,6-Tribromophenol	15.775	330	1315	0.191	ng	-0.01
15) 2-Fluorobiphenyl	12.895	172	10199	0.187	ng	-0.01
27) Fluoranthene-d10	19.063	212	17604	0.214	ng	0.00
31) Terphenyl-d14	19.662	244	13435	0.191	ng	0.00
Target Compounds						
34) Bis(2-ethylhexyl)phtha...	21.130	149	6168	0.083	ng	98
37) Benzo(b)fluoranthene	22.782	252	2368	0.021	ng	# 1

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

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