

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062124\
 Data File : BN032155.D
 Acq On : 22 Jun 2024 06:11
 Operator : MA/JU
 Sample : P2892-01
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BP-VPB-196-GW-510-512

Quant Time: Jun 22 06:49:24 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	7.639	152	11761	0.400	ng	0.00
7) Naphthalene-d8	10.410	136	39571	0.400	ng	-0.01
13) Acenaphthene-d10	14.274	164	24249	0.400	ng	-0.01
19) Phenanthrene-d10	17.029	188	49437	0.400	ng	0.00
29) Chrysene-d12	21.220	240	30341	0.400	ng	# 0.00
35) Perylene-d12	23.443	264	30036	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.241	112	3097	0.093	ng	0.00
5) Phenol-d6	6.823	99	4173	0.095	ng	0.00
8) Nitrobenzene-d5	8.798	82	7834	0.238	ng	0.00
11) 2-Methylnaphthalene-d10	12.009	152	17437	0.277	ng	-0.01
14) 2,4,6-Tribromophenol	15.775	330	1089	0.094	ng	-0.01
15) 2-Fluorobiphenyl	12.895	172	27804	0.302	ng	-0.01
27) Fluoranthene-d10	19.063	212	25440	0.189	ng	0.00
31) Terphenyl-d14	19.667	244	19407	0.262	ng	0.00
Target Compounds						
34) Bis(2-ethylhexyl)phtha...	21.130	149	26747	0.343	ng	97
37) Benzo(b)fluoranthene	22.777	252	3387	0.031	ng	# 1

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

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