

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100923\
 Data File : BN028165.D
 Acq On : 09 Oct 2023 17:35
 Operator : MA/JU
 Sample : 04662-02
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BP-VPB-RW10A-GW-520-522

Quant Time: Oct 09 18:08:39 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN100223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 09 12:40:55 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.929	152	3567	0.400	ng	0.00
7) Naphthalene-d8	10.746	136	10262	0.400	ng	0.01
13) Acenaphthene-d10	14.566	164	5373	0.400	ng	0.00
19) Phenanthrene-d10	17.319	188	10196	0.400	ng	#-0.01
29) Chrysene-d12	21.515	240	9291	0.400	ng	0.00
35) Perylene-d12	23.978	264	9437	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.473	112	1533	0.144	ng	0.00
5) Phenol-d6	7.098	99	1202	0.090	ng	0.01
8) Nitrobenzene-d5	9.134	82	2073	0.237	ng	0.01
11) 2-Methylnaphthalene-d10	12.324	152	3557	0.234	ng	0.00
14) 2,4,6-Tribromophenol	16.077	330	437	0.178	ng	0.00
15) 2-Fluorobiphenyl	13.187	172	5188	0.268	ng	-0.01
27) Fluoranthene-d10	19.356	212	6552	0.256	ng	0.00
31) Terphenyl-d14	19.941	244	6196	0.286	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.364	88	83	0.021	ng	# 25
20) 4,6-Dinitro-2-methylph...	16.760	198	4	0.030	ng	# 50
24) Pentachlorophenol	16.971	266	114	0.043	ng	# 79
34) Bis(2-ethylhexyl)phtha...	21.372	149	7357	0.354	ng	98

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

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