

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN051024\
 Data File : BN031483.D
 Acq On : 11 May 2024 04:28
 Operator : MA/JU
 Sample : P2319-03
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 TEST-2-T42

Quant Time: May 11 04:56:39 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN050924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 10 02:12:24 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.711	152	12281	0.400	ng	0.00
7) Naphthalene-d8	10.496	136	39113	0.400	ng	# 0.00
13) Acenaphthene-d10	14.338	164	22378	0.400	ng	-0.01
19) Phenanthrene-d10	17.078	188	51644	0.400	ng	#-0.01
29) Chrysene-d12	21.265	240	43898	0.400	ng	0.00
35) Perylene-d12	23.505	264	38986	0.400	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.299	112	3323	0.095	ng	0.00
5) Phenol-d6	6.866	99	2694	0.057	ng	0.00
8) Nitrobenzene-d5	8.852	82	7641	0.272	ng	-0.01
11) 2-Methylnaphthalene-d10	12.085	152	16039	0.254	ng	0.00
14) 2,4,6-Tribromophenol	15.825	330	2085	0.246	ng	-0.01
15) 2-Fluorobiphenyl	12.970	172	24199	0.280	ng	-0.01
27) Fluoranthene-d10	19.114	212	45268	0.316	ng	0.00
31) Terphenyl-d14	19.723	244	33986	0.316	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.240	88	2539	0.168	ng	# 20
33) Chrysene	21.300	228	3387	0.021	ng	95
34) Bis(2-ethylhexyl)phtha...	21.202	149	12305	0.124	ng	98
37) Benzo(b)fluoranthene	22.832	252	2950	0.021	ng	# 36
38) Benzo(k)fluoranthene	22.876	252	2826	0.021	ng	# 29

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

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