

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112822\
 Data File : BM037763.D
 Acq On : 28 Nov 2022 17:15
 Operator : CG/JU
 Sample : N5658-08
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ESQS6

Quant Time: Nov 28 22:29:27 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 28 22:24:47 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.110	152	216531	20.000	ng/u1	0.00
20) Naphthalene-d8	10.928	136	1004794	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.733	164	647193	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.468	188	1345445	20.000	ng/u1	0.00
79) Chrysene-d12	21.639	240	1236084	20.000	ng/u1	# 0.00
88) Perylene-d12	24.127	264	1066876	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.434	96	30886	5.916	ng/uL	0.00
4) Pyridine-d5	3.869	84	375675	22.105	ng/u1	0.00
7) Phenol-d5	7.257	99	693726	32.114	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.428	67	462289	32.080	ng/u1	0.00
11) 2-Chlorophenol-d4	7.634	132	517622	34.484	ng/u1	0.00
15) 4-Methylphenol-d8	8.816	113	556523	32.374	ng/u1	0.00
21) Nitrobenzene-d5	9.281	128	284396	36.104	ng/u1	0.00
24) 2-Nitrophenol-d4	10.004	143	294721	37.971	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.545	165	507434	35.725	ng/u1	0.00
31) 4-Chloroaniline-d4	11.063	131	486308	20.633	ng/u1	0.00
46) Dimethylphthalate-d6	14.139	166	1587365	34.627	ng/u1	0.00
49) Acenaphthylene-d8	14.427	160	1910191	35.302	ng/u1	0.00
54) 4-Nitrophenol-d4	14.916	143	291081	28.827	ng/u1	0.00
60) Fluorene-d10	15.721	176	1426886	35.540	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.827	200	192989	26.664	ng/u1	0.00
73) Anthracene-d10	17.568	188	2211620	35.151	ng/u1	0.00
81) Pyrene-d10	19.851	212	2469062	39.307	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.968	264	1939522	35.319	ng/u1	0.00

Target Compounds Qvalue

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

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