

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081222\
 Data File : BM036241.D
 Acq On : 12 Aug 2022 22:24
 Operator : CG/JU
 Sample : PB146838BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK838

Quant Time: Aug 12 23:55:45 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM072522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Aug 12 11:23:11 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.816	152	204577	20.000	ng/u1	0.00
20) Naphthalene-d8	10.616	136	850949	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.457	164	474122	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.198	188	939671	20.000	ng/u1	0.00
79) Chrysene-d12	21.380	240	954473	20.000	ng/u1	0.00
88) Perylene-d12	23.715	264	1056147	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.216	96	33082	6.085	ng/uL	0.00
4) Pyridine-d5	3.640	84	438579	29.447	ng/u1	0.00
7) Phenol-d5	7.004	99	530822	30.941	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.151	67	328184	28.696	ng/u1	0.00
11) 2-Chlorophenol-d4	7.351	132	448641	31.505	ng/u1	0.00
15) 4-Methylphenol-d8	8.545	113	401808	28.911	ng/u1	0.00
21) Nitrobenzene-d5	8.987	128	208215	30.879	ng/u1	0.00
24) 2-Nitrophenol-d4	9.704	143	218932	33.325	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.251	165	364787	30.701	ng/u1	0.00
31) 4-Chloroaniline-d4	10.769	131	565069	28.093	ng/u1	0.00
46) Dimethylphthalate-d6	13.880	166	1029249	31.869	ng/u1	0.00
49) Acenaphthylene-d8	14.151	160	1308074	31.744	ng/u1	0.00
54) 4-Nitrophenol-d4	14.692	143	170347	25.686	ng/u1	0.00
60) Fluorene-d10	15.451	176	917984	31.455	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.580	200	131991	26.132	ng/u1	0.00
73) Anthracene-d10	17.298	188	1399636	32.946	ng/u1	0.00
81) Pyrene-d10	19.592	212	1630173	30.817	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.568	264	1770064	33.012	ng/u1	0.00

Target Compounds Qvalue

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

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