

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082222\
 Data File : BM036392.D
 Acq On : 23 Aug 2022 16:22
 Operator : CG/JU
 Sample : PB147163BL
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK163

Quant Time: Aug 24 01:00:28 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM081622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 23 00:55:01 2022
 Response via : Initial Calibration

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

Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.792	152	227626	20.000 ng/u1	0.00
20) Naphthalene-d8	10.592	136	1014042	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.433	164	575431	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.180	188	997467	20.000 ng/u1	0.00
79) Chrysene-d12	21.362	240	728663	20.000 ng/u1	0.00
88) Perylene-d12	23.691	264	731917	20.000 ng/u1	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.204	96	37758	6.322 ng/uL	0.00
4) Pyridine-d5	3.628	84	548557	33.305 ng/u1	0.00
7) Phenol-d5	6.969	99	682951	35.150 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	469720	38.407 ng/u1	0.00
11) 2-Chlorophenol-d4	7.322	132	543585	35.607 ng/u1	0.00
15) 4-Methylphenol-d8	8.516	113	539035	37.156 ng/u1	0.00
21) Nitrobenzene-d5	8.963	128	277818	36.135 ng/u1	0.00
24) 2-Nitrophenol-d4	9.680	143	277215	35.547 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.222	165	469472	33.691 ng/u1	0.00
31) 4-Chloroaniline-d4	10.745	131	775810	35.913 ng/u1	0.00
46) Dimethylphthalate-d6	13.857	166	1471415	40.419 ng/u1	0.00
49) Acenaphthylene-d8	14.133	160	1723122	37.950 ng/u1	0.00
54) 4-Nitrophenol-d4	14.668	143	185253	26.856 ng/u1	0.00
60) Fluorene-d10	15.427	176	1267318	38.525 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.562	200	138931	27.092 ng/u1	0.00
73) Anthracene-d10	17.280	188	1748064	39.413 ng/u1	0.00
81) Pyrene-d10	19.574	212	1695816	44.864 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.538	264	1417936	39.513 ng/u1	0.00

Target Compounds Qvalue

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

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