

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092122\
 Data File : BM036634.D
 Acq On : 21 Sep 2022 21:12
 Operator : CG/JU
 Sample : PB147683BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK683

Quant Time: Sep 21 22:53:54 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 21 22:42:41 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.022	152	347677	20.000	ng/u1	0.00
20) Naphthalene-d8	10.833	136	1617040	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.645	164	1078587	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.380	188	2264081	20.000	ng/u1	0.00
79) Chrysene-d12	21.545	240	1923304	20.000	ng/u1	0.00
88) Perylene-d12	23.986	264	1754805	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	42847	4.620	ng/uL	0.00
4) Pyridine-d5	3.822	84	477377	17.649	ng/u1	0.00
7) Phenol-d5	7.175	99	770361	24.794	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.345	67	482706	25.582	ng/u1	0.00
11) 2-Chlorophenol-d4	7.551	132	621692	27.793	ng/u1	0.00
15) 4-Methylphenol-d8	8.728	113	631852	27.522	ng/u1	0.00
21) Nitrobenzene-d5	9.187	128	318948	29.141	ng/u1	0.00
24) 2-Nitrophenol-d4	9.910	143	322912	32.938	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.451	165	594524	26.168	ng/u1	0.00
31) 4-Chloroaniline-d4	10.969	131	604114	15.993	ng/u1	0.00
46) Dimethylphthalate-d6	14.057	166	2060833	28.330	ng/u1	0.00
49) Acenaphthylene-d8	14.339	160	2392396	27.726	ng/u1	0.00
54) 4-Nitrophenol-d4	14.833	143	353635	25.064	ng/u1	0.00
60) Fluorene-d10	15.633	176	1813366	27.816	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	215227	24.027	ng/u1	0.00
73) Anthracene-d10	17.480	188	2838596	27.421	ng/u1	0.00
81) Pyrene-d10	19.762	212	3125578	33.033	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.827	264	2451003	28.589	ng/u1	0.00

Target Compounds Qvalue

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

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