

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101723\
 Data File : BM042329.D
 Acq On : 17 Oct 2023 20:14
 Operator : MA/JU
 Sample : PB156422BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK422

Quant Time: Oct 18 02:16:47 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM101223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 18 02:04:27 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.010	152	114998	20.000	ng/u1	0.00
20) Naphthalene-d8	10.833	136	518183	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.651	164	313147	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.404	188	661876	20.000	ng/u1	0.00
79) Chrysene-d12	21.580	240	563458	20.000	ng/u1	0.00
88) Perylene-d12	24.056	264	587673	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	28366	7.707	ng/uL	0.00
4) Pyridine-d5	3.804	84	415194	38.090	ng/u1	0.00
7) Phenol-d5	7.157	99	482664	37.872	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.351	67	340214	39.507	ng/u1	0.00
11) 2-Chlorophenol-d4	7.534	132	340822	38.874	ng/u1	0.00
15) 4-Methylphenol-d8	8.716	113	368798	36.540	ng/u1	0.00
21) Nitrobenzene-d5	9.210	128	170463	38.734	ng/u1	0.00
24) 2-Nitrophenol-d4	9.922	143	182396	38.771	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.445	165	311569	36.291	ng/u1	0.00
31) 4-Chloroaniline-d4	10.998	131	494965	36.617	ng/u1	0.00
46) Dimethylphthalate-d6	14.068	166	1097053	40.414	ng/u1	0.00
49) Acenaphthylene-d8	14.357	160	1188179	39.992	ng/u1	0.00
54) 4-Nitrophenol-d4	14.863	143	186540	34.165	ng/u1	0.00
60) Fluorene-d10	15.645	176	908216	40.700	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.774	200	141409	30.499	ng/u1	0.00
73) Anthracene-d10	17.504	188	1441767	42.523	ng/u1	0.00
81) Pyrene-d10	19.798	212	1657410	48.263	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.897	264	1472965	45.664	ng/u1	0.00

Target Compounds Qvalue

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

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