

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111723\
 Data File : BM042869.D
 Acq On : 18 Nov 2023 07:02
 Operator : MA/JU
 Sample : PB156976BL
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Nov 18 21:06:55 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 15 22:28:07 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.816	152	25543	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.628	136	100897	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.474	164	62726	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.227	188	134332	20.000	ng/u1	-0.01
79) Chrysene-d12	21.415	240	131202	20.000	ng/u1	0.00
88) Perylene-d12	23.768	264	166066	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	5072	6.231	ng/uL	0.00
4) Pyridine-d5	3.610	84	57915	28.602	ng/u1	0.00
7) Phenol-d5	6.987	99	63284	25.456	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.163	67	53438	30.591	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.340	132	50704	26.433	ng/u1	-0.01
15) 4-Methylphenol-d8	8.540	113	46829	23.924	ng/u1	-0.01
21) Nitrobenzene-d5	9.098	128	281	0.318	ng/u1	0.07
24) 2-Nitrophenol-d4	9.734	143	23443	22.654	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.257	165	39918	20.745	ng/u1	0.00
31) 4-Chloroaniline-d4	10.810	131	52518	21.496	ng/u1	0.00
46) Dimethylphthalate-d6	13.898	166	159204	26.817	ng/u1	0.00
49) Acenaphthylene-d8	14.174	160	171192	26.951	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.822	143	104	0.155	ng/u1	0.08
60) Fluorene-d10	15.469	176	135630	27.590	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.621	200	15755	15.484	ng/u1	0.00
73) Anthracene-d10	17.327	188	202971	26.961	ng/u1	-0.01
81) Pyrene-d10	19.621	212	277159	32.259	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.615	264	285688	28.760	ng/u1	-0.01

Target Compounds Qvalue

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

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