

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110723\
 Data File : BP017948.D
 Acq On : 08 Nov 2023 05:02
 Operator : MA/JU
 Sample : PB156700BL
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK700

Quant Time: Nov 08 05:32:45 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110623.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 07 03:40:56 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	75930	20.000	ng/u1	0.00
20) Naphthalene-d8	10.669	136	315775	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.516	164	197325	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.292	188	432267	20.000	ng/u1	0.00
79) Chrysene-d12	21.410	240	389197	20.000	ng/u1	0.00
88) Perylene-d12	23.892	264	431017	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	12165	6.132	ng/uL	0.00
4) Pyridine-d5	3.675	84	157828	28.665	ng/u1	0.00
7) Phenol-d5	7.022	99	196340	27.574	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.204	67	129334	30.111	ng/u1	0.00
11) 2-Chlorophenol-d4	7.381	132	160621	29.357	ng/u1	0.00
15) 4-Methylphenol-d8	8.575	113	160333	27.243	ng/u1	0.00
21) Nitrobenzene-d5	9.057	128	79518	31.511	ng/u1	0.00
24) 2-Nitrophenol-d4	9.769	143	88921	30.594	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.298	165	151265	28.388	ng/u1	0.00
31) 4-Chloroaniline-d4	10.845	131	207356	26.854	ng/u1	0.00
46) Dimethylphthalate-d6	13.939	166	520997	30.111	ng/u1	0.00
49) Acenaphthylene-d8	14.216	160	561997	29.601	ng/u1	0.00
54) 4-Nitrophenol-d4	14.775	143	59488	22.361	ng/u1	0.00
60) Fluorene-d10	15.516	176	427620	29.528	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.669	200	70710	22.369	ng/u1	0.00
73) Anthracene-d10	17.392	188	674594	30.015	ng/u1	0.00
81) Pyrene-d10	19.639	212	783206	31.701	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.721	264	739849	30.721	ng/u1	-0.01

Target Compounds Qvalue

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

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