

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110723\
 Data File : BP017947.D
 Acq On : 08 Nov 2023 04:28
 Operator : MA/JU
 Sample : PB156698BL
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK698

Quant Time: Nov 08 05:02:10 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	7.858	152	75155	20.000	ng/u1	0.00
20) Naphthalene-d8	10.669	136	318966	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.516	164	197262	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.292	188	445023	20.000	ng/u1	0.00
79) Chrysene-d12	21.410	240	385442	20.000	ng/u1	0.00
88) Perylene-d12	23.892	264	418389	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	13850	7.053	ng/uL	0.00
4) Pyridine-d5	3.675	84	190211	34.903	ng/u1	0.00
7) Phenol-d5	7.022	99	228295	32.393	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.205	67	150490	35.398	ng/u1	0.00
11) 2-Chlorophenol-d4	7.381	132	187738	34.667	ng/u1	0.00
15) 4-Methylphenol-d8	8.575	113	189505	32.532	ng/u1	0.00
21) Nitrobenzene-d5	9.057	128	95837	37.598	ng/u1	0.00
24) 2-Nitrophenol-d4	9.769	143	107691	36.681	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.299	165	179292	33.312	ng/u1	0.00
31) 4-Chloroaniline-d4	10.846	131	245196	31.437	ng/u1	0.00
46) Dimethylphthalate-d6	13.940	166	619220	35.799	ng/u1	0.00
49) Acenaphthylene-d8	14.216	160	669058	35.251	ng/u1	0.00
54) 4-Nitrophenol-d4	14.775	143	72253	27.168	ng/u1	0.00
60) Fluorene-d10	15.516	176	506137	34.961	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.669	200	85941	26.407	ng/u1	0.00
73) Anthracene-d10	17.392	188	806169	34.841	ng/u1	0.00
81) Pyrene-d10	19.639	212	938602	38.361	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.727	264	860510	36.809	ng/u1	0.00

Target Compounds Qvalue

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

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