

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110723\
 Data File : BP017962.D
 Acq On : 08 Nov 2023 14:15
 Operator : MA/JU
 Sample : PB156829BL
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK829

Quant Time: Nov 08 22:17:21 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	7.852	152	87902	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.669	136	362660	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.516	164	236129	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.292	188	531851	20.000	ng/u1	0.00
79) Chrysene-d12	21.410	240	490085	20.000	ng/u1	0.00
88) Perylene-d12	23.892	264	532080	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	11580	5.042	ng/uL	0.00
4) Pyridine-d5	3.676	84	151516	23.771	ng/u1	0.00
7) Phenol-d5	7.022	99	176241	21.380	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.205	67	115895	23.307	ng/u1	0.00
11) 2-Chlorophenol-d4	7.375	132	145781	23.015	ng/u1	-0.01
15) 4-Methylphenol-d8	8.575	113	146157	21.452	ng/u1	0.00
21) Nitrobenzene-d5	9.058	128	70897	24.463	ng/u1	0.00
24) 2-Nitrophenol-d4	9.769	143	79276	23.749	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.299	165	137069	22.399	ng/u1	0.00
31) 4-Chloroaniline-d4	10.846	131	187910	21.190	ng/u1	0.00
46) Dimethylphthalate-d6	13.940	166	501147	24.204	ng/u1	0.00
49) Acenaphthylene-d8	14.216	160	526454	23.172	ng/u1	0.00
54) 4-Nitrophenol-d4	14.775	143	52379	16.453	ng/u1	0.00
60) Fluorene-d10	15.516	176	408797	23.589	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.669	200	67847	17.444	ng/u1	0.00
73) Anthracene-d10	17.392	188	660213	23.875	ng/u1	0.00
81) Pyrene-d10	19.639	212	784982	25.232	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.727	264	728612	24.508	ng/u1	0.00

Target Compounds	Qvalue

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

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