

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100423\
 Data File : BG059275.D
 Acq On : 4 Oct 2023 10:53
 Operator : MA/JU
 Sample : PB155829BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK829

Quant Time: Oct 04 22:58:28 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 28 04:33:25 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.235	152	52010	20.000	ng/u1	0.00
20) Naphthalene-d8	11.072	136	242534	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.862	164	153193	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.612	188	336020	20.000	ng/u1	0.00
79) Chrysene-d12	21.913	240	311010	20.000	ng/u1	0.00
88) Perylene-d12	25.391	264	374397	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.528	96	8175	5.932	ng/uL	0.00
4) Pyridine-d5	3.963	84	122252	30.405	ng/u1	0.00
7) Phenol-d5	7.353	99	159582	30.653	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.547	67	99975	30.061	ng/u1	0.00
11) 2-Chlorophenol-d4	7.747	132	113062	32.307	ng/u1	0.00
15) 4-Methylphenol-d8	8.922	113	121876	30.285	ng/u1	0.00
21) Nitrobenzene-d5	9.427	128	57049	33.333	ng/u1	0.00
24) 2-Nitrophenol-d4	10.144	143	64813	38.576	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.673	165	114215	33.092	ng/u1	0.00
31) 4-Chloroaniline-d4	11.219	131	174135	31.184	ng/u1	0.00
46) Dimethylphthalate-d6	14.263	166	379970	34.306	ng/u1	0.00
49) Acenaphthylene-d8	14.568	160	450937	32.436	ng/u1	0.00
54) 4-Nitrophenol-d4	15.056	143	57260	29.605	ng/u1	0.00
60) Fluorene-d10	15.849	176	340478	32.298	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.961	200	46786	27.088	ng/u1	0.00
73) Anthracene-d10	17.712	188	526406	34.381	ng/u1	0.00
81) Pyrene-d10	19.986	212	587650	33.740	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.150	264	610709	33.944	ng/u1	-0.01

Target Compounds Qvalue

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

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