

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092622\
 Data File : BM036768.D
 Acq On : 27 Sep 2022 05:40
 Operator : CG/JU
 Sample : PB147898BL
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK898

Quant Time: Sep 27 06:25:52 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Sep 26 23:10:22 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.998	152	406986	20.000	ng/u1	0.00
20) Naphthalene-d8	10.810	136	1803876	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.621	164	1136523	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.363	188	2182406	20.000	ng/u1	0.00
79) Chrysene-d12	21.527	240	1608876	20.000	ng/u1	0.00
88) Perylene-d12	23.956	264	1393618	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	65227	6.008	ng/uL	0.00
4) Pyridine-d5	3.810	84	833173	26.314	ng/u1	0.00
7) Phenol-d5	7.157	99	1164612	32.021	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.328	67	729977	33.049	ng/u1	0.00
11) 2-Chlorophenol-d4	7.528	132	920739	35.163	ng/u1	0.00
15) 4-Methylphenol-d8	8.704	113	930164	34.612	ng/u1	-0.01
21) Nitrobenzene-d5	9.163	128	463680	37.977	ng/u1	0.00
24) 2-Nitrophenol-d4	9.886	143	455848	41.682	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.428	165	835821	32.978	ng/u1	0.00
31) 4-Chloroaniline-d4	10.945	131	1246998	29.592	ng/u1	0.00
46) Dimethylphthalate-d6	14.039	166	2620618	34.188	ng/u1	0.00
49) Acenaphthylene-d8	14.322	160	3136420	34.496	ng/u1	0.00
54) 4-Nitrophenol-d4	14.816	143	395824	26.624	ng/u1	0.00
60) Fluorene-d10	15.610	176	2330878	33.932	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.727	200	241115	27.924	ng/u1	0.00
73) Anthracene-d10	17.462	188	3452770	34.602	ng/u1	0.00
81) Pyrene-d10	19.745	212	3587503	45.325	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.797	264	2399512	35.242	ng/u1	0.00

Target Compounds Qvalue

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

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