

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091023\
 Data File : BP017092.D
 Acq On : 11 Sep 2023 06:28
 Operator : MA/JU
 Sample : PB155340BL
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK340

Quant Time: Sep 11 07:18:30 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Sep 11 06:13:29 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.958	152	744855	20.000	ng/u1	0.00
20) Naphthalene-d8	10.787	136	3388523	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.622	164	2200954	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.386	188	4855154	20.000	ng/u1	0.00
79) Chrysene-d12	21.486	240	4394355	20.000	ng/u1	0.00
88) Perylene-d12	24.021	264	4757287	20.000	ng/u1	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.317	96	111529	5.715	ng/uL	0.00
4) Pyridine-d5	3.752	84	1636644	27.811	ng/u1	0.00
7) Phenol-d5	7.111	99	2040016	30.130	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.305	67	1372321	30.220	ng/u1	0.00
11) 2-Chlorophenol-d4	7.481	132	1523966	31.142	ng/u1	0.00
15) 4-Methylphenol-d8	8.669	113	1630484	30.442	ng/u1	0.00
21) Nitrobenzene-d5	9.163	128	792053	32.832	ng/u1	0.00
24) 2-Nitrophenol-d4	9.875	143	877867	35.258	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.399	165	1460290	30.930	ng/u1	0.00
31) 4-Chloroaniline-d4	10.946	131	2220708	29.333	ng/u1	0.00
46) Dimethylphthalate-d6	14.034	166	5047092	32.227	ng/u1	0.00
49) Acenaphthylene-d8	14.322	160	5529652	31.008	ng/u1	0.00
54) 4-Nitrophenol-d4	14.840	143	962920	29.803	ng/u1	0.00
60) Fluorene-d10	15.616	176	4192133	31.544	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	783501	28.596	ng/u1	0.00
73) Anthracene-d10	17.492	188	6616870	31.795	ng/u1	0.00
81) Pyrene-d10	19.727	212	7593554	32.814	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.857	264	7352508	32.194	ng/u1	0.00

Target Compounds	Qvalue

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

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