

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091224\
 Data File : BP021895.D
 Acq On : 12 Sep 2024 14:11
 Operator : RC/JU
 Sample : PB163275BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK275

Quant Time: Sep 13 02:41:55 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 12 12:48:41 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	83412	20.000	ng/u1	0.03
20) Naphthalene-d8	10.528	136	324358	20.000	ng/u1	# 0.02
38) Acenaphthene-d10	14.363	164	240897	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.157	188	555943	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.586	240	607517	20.000	ng/u1	# 0.00
88) Perylene-d12	24.909	264	726406	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	10628	4.122	ng/uL	0.01
4) Pyridine-d5	3.681	84	135722	17.877	ng/u1	0.02
7) Phenol-d5	6.934	99	161668	22.683	ng/u1	0.02
9) Bis-(2-Chloroethyl)eth...	7.093	67	100072	24.174	ng/u1	0.02
11) 2-Chlorophenol-d4	7.299	132	137742	27.763	ng/u1	0.02
15) 4-Methylphenol-d8	8.457	113	133647	23.486	ng/u1	0.03
21) Nitrobenzene-d5	8.899	128	67571	33.449	ng/u1	0.02
24) 2-Nitrophenol-d4	9.616	143	73194	31.873	ng/u1	0.02
28) 2,4-Dichlorophenol-d3	10.151	165	142052	27.933	ng/u1	0.02
31) 4-Chloroaniline-d4	10.657	131	180210	24.315	ng/u1	0.02
46) Dimethylphthalate-d6	13.769	166	524519	28.982	ng/u1	0.00
49) Acenaphthylene-d8	14.057	160	531329	26.764	ng/u1	0.00
54) 4-Nitrophenol-d4	14.563	143	82743	24.185	ng/u1	0.00
60) Fluorene-d10	15.363	176	436910	25.398	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.486	200	81512	25.628	ng/u1	0.00
73) Anthracene-d10	17.257	188	736824	28.532	ng/u1	0.00
81) Pyrene-d10	19.622	212	943715	28.077	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.686	264	1042164	27.568	ng/u1	0.00

Target Compounds Qvalue

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

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